

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012142.D
 Acq On : 14 Oct 2022 17:35
 Operator : CG/JU
 Sample : N4982-16
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 14 23:04:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	259022	20.000	ng/ul	0.00
20) Naphthalene-d8	10.652	136	1050018	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.493	164	566241	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	1207065	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1346375	20.000	ng/ul	0.00
88) Perylene-d12	23.763	264	1426336	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	27598	4.201	ng/uL	0.00
4) Pyridine-d5	3.699	84	160020	9.159	ng/ul	0.00
7) Phenol-d5	7.016	99	171486	8.232	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	397510	33.617	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.375	132	449684	27.194	ng/ul	0.00
15) 4-Methylphenol-d8	8.558	113	276263	17.265	ng/ul	0.00
21) Nitrobenzene-d5	9.010	128	261542	35.268	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.734	143	259517	34.513	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.275	165	422753	29.009	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	434065	20.212	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1214747	33.874	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	1495091	33.715	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	36818	5.230	ng/ul	0.00
60) Fluorene-d10	15.493	176	1121617	34.570	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	151781	23.814	ng/ul	0.00
73) Anthracene-d10	17.351	188	1862163	35.527	ng/ul	0.00
81) Pyrene-d10	19.592	212	2364764	34.284	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	2474781	35.692	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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